```
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TERMINAL (ENTER 1, 2, 3, OR ?):2


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 NEWS  1           Web Page URLs for STN Seminar Schedule - N. America
 NEWS  2           "Ask CAS" for self-help around the clock
 NEWS  3  SEP 01   New pricing for the Save Answers for SciFinder Wizard within
                   STN Express with Discover!
 NEWS  4  OCT 28   KOREAPAT now available on STN
 NEWS  5  NOV 30   PHAR reloaded with additional data
 NEWS  6  DEC 01   LISA now available on STN
 NEWS  7  DEC 09   12 databases to be removed from STN on December 31, 2004
 NEWS  8  DEC 15   MEDLINE update schedule for December 2004
 NEWS  9  DEC 17   ELCOM reloaded; updating to resume; current-awareness
                   alerts (SDIs) affected
 NEWS  10 DEC 17   COMPUAB reloaded; updating to resume; current-awareness
                   alerts (SDIs) affected
 NEWS  11 DEC 17   SOLIDSTATE reloaded; updating to resume; current-awareness
                   alerts (SDIs) affected
 NEWS  12 DEC 17   CERAB reloaded; updating to resume; current-awareness
                   alerts (SDIs) affected
 NEWS  13 DEC 17   THREE NEW FIELDS ADDED TO IFIPAT/IFIUDB/IFICDB
 NEWS  14 DEC 30   EPFULL: New patent full text database to be available on STN
 NEWS  15 DEC 30   CAPLUS - PATENT COVERAGE EXPANDED
 NEWS  16 JAN 03   No connect-hour charges in EPFULL during January and
                   February 2005
 NEWS  17 FEB 25   CA/CAPLUS - Russian Agency for Patents and Trademarks
                   (ROSPATENT) added to list of core patent offices covered
 NEWS  18 FEB 10   STN Patent Forums to be held in March 2005
 NEWS  19 FEB 16   STN User Update to be held in conjunction with the 229th ACS
                   National Meeting on March 13, 2005
 NEWS  20 FEB 28   PATDPAFULL - New display fields provide for legal status
                   data from INPADOC
 NEWS  21 FEB 28   BABS - Current-awareness alerts (SDIs) available
 NEWS  22 FEB 28   MEDLINE/LMEDLINE reloaded

 NEWS EXPRESS  JANUARY 10 CURRENT WINDOWS VERSION IS V7.01a, CURRENT
               MACINTOSH VERSION IS V6.0c(ENG) AND V6.0Jc(JP),
               AND CURRENT DISCOVER FILE IS DATED 10 JANUARY 2005

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  * * * * * * * * * * * * * * * STN  Columbus  * * * * * * * * * * * * * * * *
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FILE 'HOME' ENTERED AT 08:17:57 ON 02 MAR 2005

=> FIL STNGUIDE
COST IN U.S. DOLLARS

	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	0.21	0.21

FILE 'STNGUIDE' ENTERED AT 08:18:02 ON 02 MAR 2005
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AND TECHNOLOGY CORPORATION, AND FACHINFORMATIONSZENTRUM KARLSRUHE

FILE CONTAINS CURRENT INFORMATION.
LAST RELOADED: Feb 25, 2005 (20050225/UP).

=> FIL HOME
COST IN U.S. DOLLARS

	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	0.06	0.27

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=> fil reg
COST IN U.S. DOLLARS

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FULL ESTIMATED COST	0.21	0.48

FILE 'REGISTRY' ENTERED AT 08:18:15 ON 02 MAR 2005
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STRUCTURE FILE UPDATES: 28 FEB 2005 HIGHEST RN 839671-97-5
DICTIONARY FILE UPDATES: 28 FEB 2005 HIGHEST RN 839671-97-5

TSCA INFORMATION NOW CURRENT THROUGH JANUARY 18, 2005

Please note that search-term pricing does apply when
conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. For more
information enter HELP PROP at an arrow prompt in the file or refer
to the file summary sheet on the web at:
http://www.cas.org/ONLINE/DBSS/registryss.html

=>
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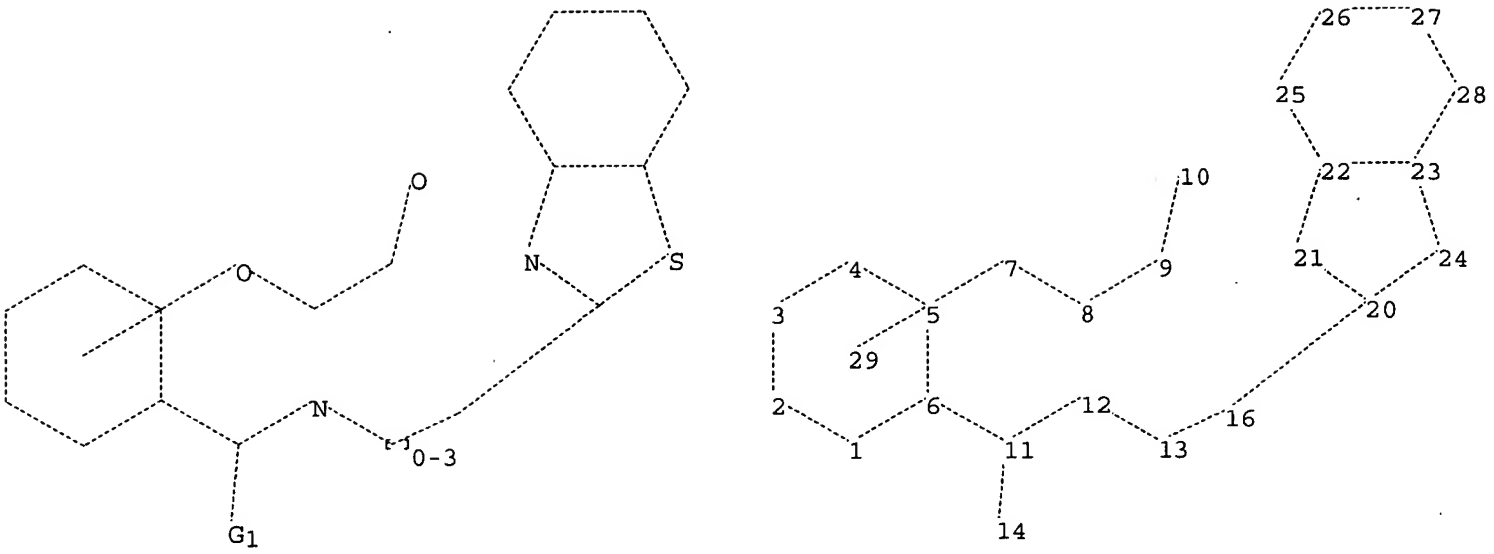

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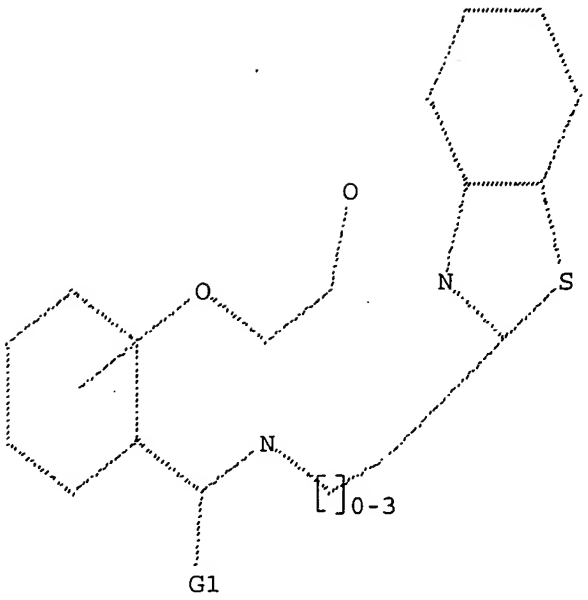
7 8 9 10 11 12 13 14 16
ring nodes :
1 2 3 4 5 6 20 21 22 23 24 25 26 27 28
chain bonds :
6-11 7-8 8-9 9-10 11-12 11-14 12-13 13-16 16-20
ring bonds :
1-2 1-6 2-3 3-4 4-5 5-6 20-21 20-24 21-22 22-23 22-25 23-24 23-28 25-26 26-27 27-28
exact/norm bonds :
1-2 1-6 2-3 3-4 4-5 5-6 6-11 7-8 8-9 9-10 11-12 11-14 12-13 13-16 16-20 20-21 20-24 21-22 22-23 22-25 23-24 23-28 25-26 26-27 27-28
isolated ring systems :
containing 1 : 20 :

G1:O,S,N

Match level :
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:CLASS 8:CLASS 9:CLASS 10:CLASS 11:CLASS 12:CLASS 13:CLASS 14:CLASS 16:CLASS 20:Atom 21:Atom 22:Atom 23:Atom 24:Atom 25:Atom 26:Atom 27:Atom 28:Atom 29:CLASS

L1 STRUCTURE UPLOADED

=> d
L1 HAS NO ANSWERS
L1 STR



G1 O,S,N

Structure attributes must be viewed using STN Express query preparation.

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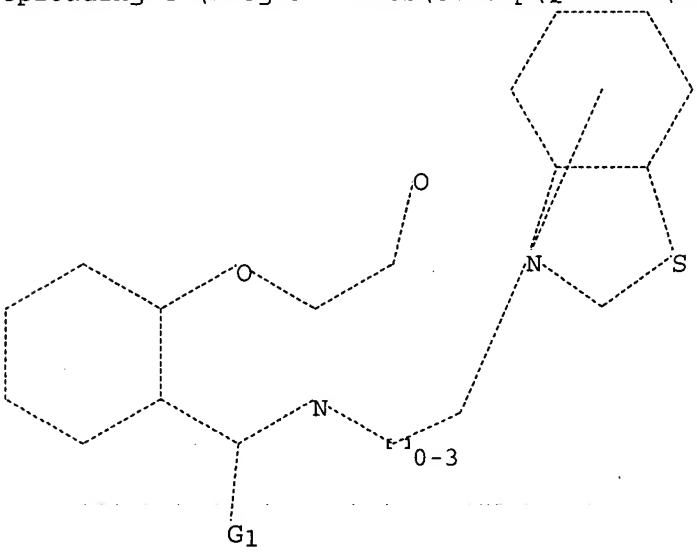

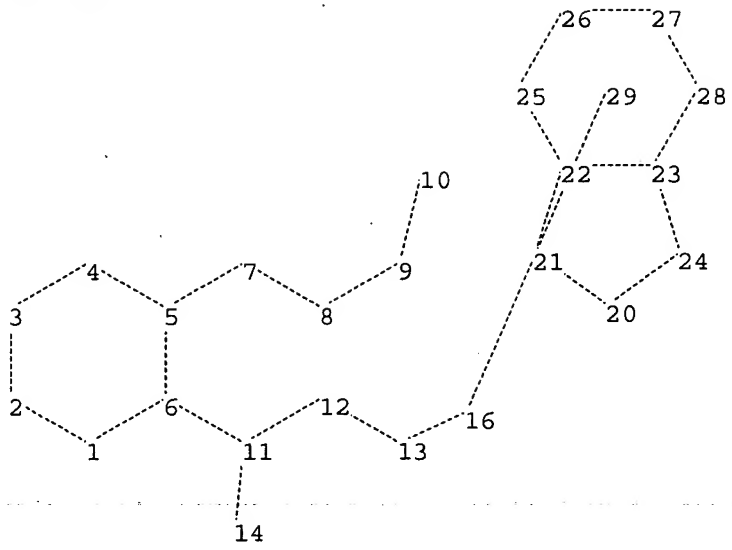

chain nodes :
7 8 9 10 11 12 13 14 16
ring nodes :
1 2 3 4 5 6 20 21 22 23 24 25 26 27 28
chain bonds :
5-7 6-11 7-8 8-9 9-10 11-12 11-14 12-13 13-16
ring bonds :
1-2 1-6 2-3 3-4 4-5 5-6 20-21 20-24 21-22 22-23 22-25 23-24 23-28 25-26 26-27
27-28
exact/norm bonds :
1-2 1-6 2-3 3-4 4-5 5-6 5-7 6-11 7-8 8-9 9-10 11-12 11-14 12-13 13-16 20-21
20-24 21-22 22-23 22-25 23-24 23-28 25-26 26-27 27-28
isolated ring systems :
containing 1 : 20 :

G1:O,S,N

Match level :
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:CLASS 8:CLASS 9:CLASS 10:CLASS 11:CLASS
12:CLASS 13:CLASS 14:CLASS 16:CLASS 20:Atom 21:Atom 22:Atom 23:Atom 24:Atom 25:Atom
26:Atom 27:Atom 28:Atom 29:CLASS

L2 STRUCTURE UPLOADED

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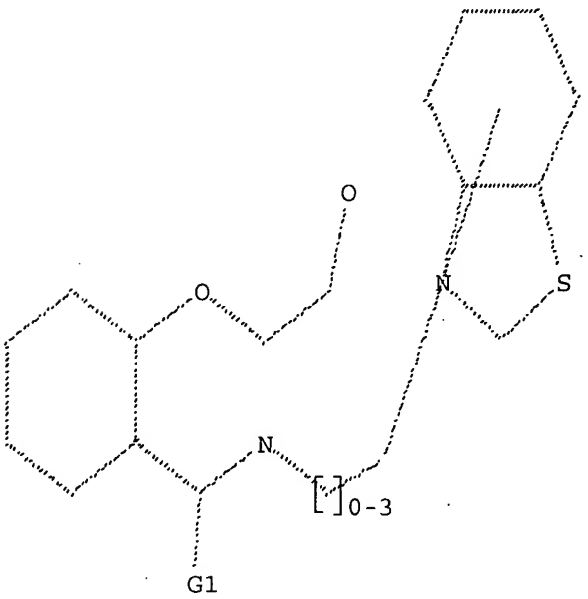
L2 HAS NO ANSWERS
L2 STR



G1 O,S,N

Structure attributes must be viewed using STN Express query preparation.

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SAMPLE SCREEN SEARCH COMPLETED - 0 TO ITERATE

100.0% PROCESSED 0 ITERATIONS 0 ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 0 TO 0
PROJECTED ANSWERS: 0 TO 0

L3 0 SEA SSS SAM L1

=> s l1 full
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FULL SCREEN SEARCH COMPLETED - 21 TO ITERATE

100.0% PROCESSED 21 ITERATIONS 8 ANSWERS
SEARCH TIME: 00.00.01

L4 8 SEA SSS FUL L1

=> s l2
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SAMPLE SCREEN SEARCH COMPLETED - 2 TO ITERATE

100.0% PROCESSED 2 ITERATIONS 0 ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 2 TO 124
PROJECTED ANSWERS: 0 TO 0

L5 0 SEA SSS SAM L2

=> s l2 full
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100.0% PROCESSED 39 ITERATIONS 0 ANSWERS
SEARCH TIME: 00.00.01

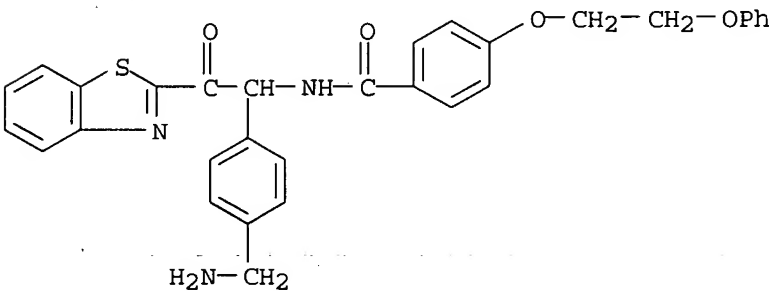
L6 0 SEA SSS FUL L2

=> s l4 and caplus/lc
44163940 CAPLUS/LC
L7 7 L4 AND CAPLUS/LC

=> s l4 not l7
L8 1 L4 NOT L7

=> d

L8 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2005 ACS on STN
RN 334989-04-7 REGISTRY
CN Benzamide, N-[1-[4-(aminomethyl)phenyl]-2-(2-benzothiazolyl)-2-oxoethyl]-4-(2-phenoxyethoxy)- (9CI) (CA INDEX NAME)
FS 3D CONCORD
MF C31 H27 N3 O4 S
CI COM
SR CA



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

=> fil caplus
COST IN U.S. DOLLARS SINCE FILE ENTRY TOTAL SESSION
FULL ESTIMATED COST 329.96 330.44

FILE 'CAPLUS' ENTERED AT 08:20:51 ON 02 MAR 2005
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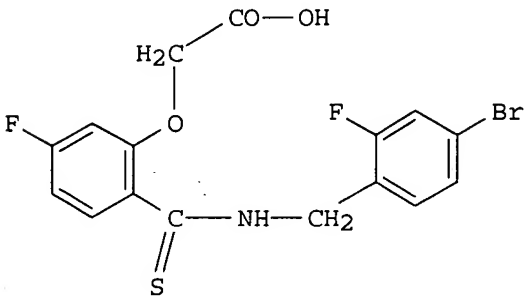
FILE COVERS 1907 - 2 Mar 2005 VOL 142 ISS 10
FILE LAST UPDATED: 1 Mar 2005 (20050301/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> s l7
L9 3 L7

=> d ibib abs hitstr l9 1-3

L9 ANSWER 1 OF 3 CAPLUS COPYRIGHT 2005 ACS on STN
ACCESSION NUMBER: 2004:819180 CAPLUS
DOCUMENT NUMBER: 141:374428
TITLE: Design and synthesis of highly potent and selective (2-arylcarbamoyl-phenoxy)-acetic acid inhibitors of aldose reductase for treatment of chronic diabetic complications
AUTHOR(S): Van Zandt, Michael C.; Sibley, Evelyn O.; McCann, Erin E.; Combs, Kerry J.; Flam, Brenda; Sawicki, Diane R.; Sabetta, Al; Carrington, Anne; Sredy, Janet; Howard, Eduardo; Mitschler, Andre; Podjarny, Alberto D.
CORPORATE SOURCE: The Institute for Diabetes Discovery, LLC, Branford, CT, 06405, USA
SOURCE: Bioorganic & Medicinal Chemistry (2004), 12(21), 5661-5675
CODEN: BMECEP; ISSN: 0968-0896
PUBLISHER: Elsevier Ltd.
DOCUMENT TYPE: Journal
LANGUAGE: English
GI



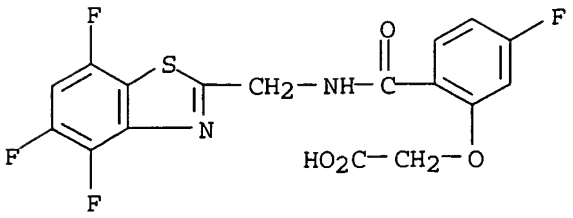
I

AB Recent efforts to identify treatments for chronic diabetic complications have resulted in the discovery of a novel series of highly potent and selective (2-arylcarbamoyl-phenoxy)-acetic acid aldose reductase inhibitors. The compound class features a core template that utilizes an intramol. hydrogen bond to position the key structural elements of the pharmacophore in a conformation, which promotes a high binding affinity. The lead candidate, I, 5-fluoro-2-(4-bromo-2-fluoro-benzylthiocarbamoyl)-phenoxyacetic acid, inhibits aldose reductase with an IC50 of 30 nM, while being 1100 times less active against aldehyde reductase, a related enzyme involved in the detoxification of reactive aldehydes. In addition, I lowers nerve sorbitol levels with an ED50 of 31 mg/kg/d po in the 4-day STZ-induced diabetic rat model.

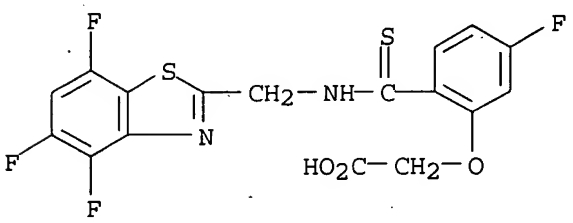
IT **314297-79-5P 314297-80-8P 314297-81-9P 314297-82-0P**

RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(design and synthesis of (arylcarbamoylphenoxy)acetic acid inhibitors of aldose reductase for treatment of chronic diabetic complications)

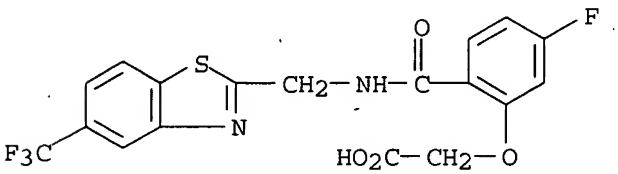
RN 314297-79-5 CAPLUS
CN Acetic acid, [5-fluoro-2-[[[(4,5,7-trifluoro-2-benzothiazolyl)methyl]amino]carbonyl]phenoxy]- (9CI) (CA INDEX NAME)



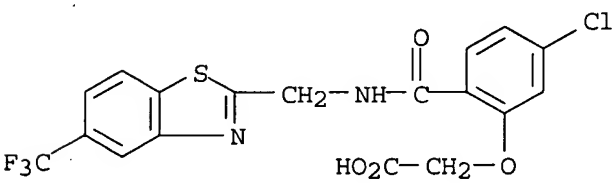
RN 314297-80-8 CAPLUS
CN Acetic acid, [5-fluoro-2-[thioxo[[(4,5,7-trifluoro-2-benzothiazolyl)methyl]amino]methyl]phenoxy]- (9CI) (CA INDEX NAME)



RN 314297-81-9 CAPLUS
CN Acetic acid, [5-fluoro-2-[[[[5-(trifluoromethyl)-2-benzothiazolyl]methyl]amino]carbonyl]phenoxy]- (9CI) (CA INDEX NAME)



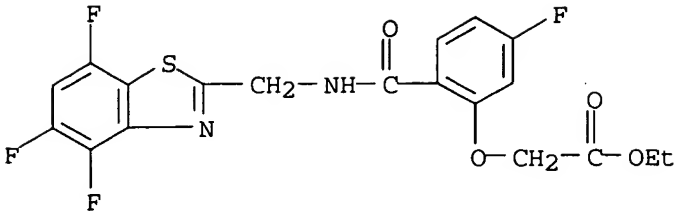
RN 314297-82-0 CAPLUS
CN Acetic acid, [5-chloro-2-[[[[5-(trifluoromethyl)-2-benzothiazolyl]methyl]amino]carbonyl]phenoxy]- (9CI) (CA INDEX NAME)



IT **314298-38-9P**
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)
(design and synthesis of (arylcarbamoylphenoxy)acetic acid inhibitors of aldose reductase for treatment of chronic diabetic complications)

RN 314298-38-9 CAPLUS
CN Acetic acid, [5-fluoro-2-[[[(4,5,7-trifluoro-2-benzothiazolyl)methyl]amino]carbonyl]phenoxy]-, ethyl ester (9CI) (CA

INDEX NAME)



REFERENCE COUNT: 44 THERE ARE 44 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L9 ANSWER 2 OF 3 CAPLUS COPYRIGHT 2005 ACS on STN
ACCESSION NUMBER: 2001:283939 CAPLUS
DOCUMENT NUMBER: 134:311433
TITLE: Preparation of (hetero)arylmethylamines as tryptase inhibitors
INVENTOR(S): Lively, Sarah Elizabeth; Waszkowycz, Bohdan; Harrison, Martin James; Clase, Juha Andrew; Naylor, Neil Jason
PATENT ASSIGNEE(S): Protherics Molecular Design Limited, UK
SOURCE: PCT Int. Appl., 106 pp.
CODEN: PIXXD2
DOCUMENT TYPE: Patent
LANGUAGE: English
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2001027096	A1	20010419	WO 2000-GB3832	20001005

W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CR, CU, CZ, DE, DK, DM, DZ, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, US, UZ, VN, YU, ZA, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM
RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG

PRIORITY APPLN. INFO.: GB 1999-23710 A 19991008
OTHER SOURCE(S): MARPAT 134:311433
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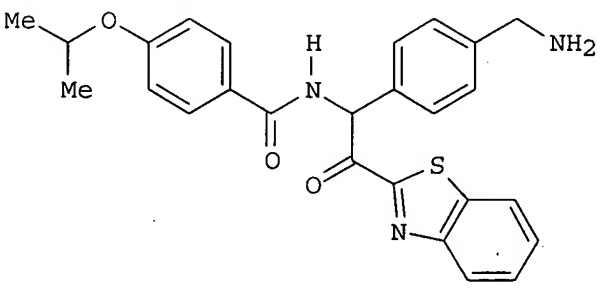

I

AB R1CHR2(CH2)aZCH2NH2 [R1 = H, NH2, NHZ1(CH2)bR3; R2 = H when R1 = NHZ1(CH2)bR3 or COR4; R3 = alk(en)yl, heterocyclyl, aryl, etc.; R4 = COR5, CF2R6, 2-(benz)oxazolyl, 2-(benz)imidazolyl, etc.; R5 = (fluoro)alkyl, alkoxy, aryl, etc.; R6 = F, (fluoro)alkyl, aryl, etc.; Z = 1,4-phenylene, 5-membered heteroarylene, etc.; Z1 = bond, CO CO2, CONH, SO2; a = 0-2; b = 0-4] were prepared as tryptase inhibitors (no data). Thus, 4-BrC6H4CH2CO2H was converted in 7 steps to 4-(BocHNH2C)C6H4CH(NH2)CO2Me which was amidated by 4-(Me2HCO)C6H4CO2H and the product condensed with

benzothiazole to give, after deprotection, title compound I.

IT **334989-05-8P**
RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(preparation of (hetero)arylmethylamines as tryptase inhibitors)

RN 334989-05-8 CAPLUS

CN Benzamide, N-[1-[4-(aminomethyl)phenyl]-2-(2-benzothiazolyl)-2-oxoethyl]-4-(2-phenoxyethoxy)-, mono(trifluoroacetate) (9CI) (CA INDEX NAME)

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CRN 334989-04-7
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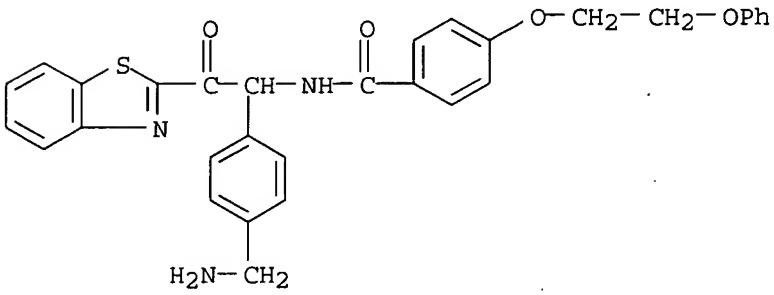

CM 2

CRN 76-05-1
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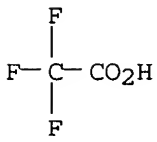


REFERENCE COUNT: 9 THERE ARE 9 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L9 ANSWER 3 OF 3 CAPLUS COPYRIGHT 2005 ACS on STN

ACCESSION NUMBER: 2001:12407 CAPLUS
DOCUMENT NUMBER: 134:71392
TITLE: Preparation and effect of Substituted phenoxyacetic acids in complications arising from diabetes mellitus
INVENTOR(S): Van Zandt, Michael C.
PATENT ASSIGNEE(S): The Institutes for Pharmaceutical Discovery, Llc, USA
SOURCE: PCT Int. Appl., 144 pp.
CODEN: PIXXD2
DOCUMENT TYPE: Patent
LANGUAGE: English
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2001000566	A2	20010104	WO 2000-US17377	20000623
WO 2001000566	A3	20020207		

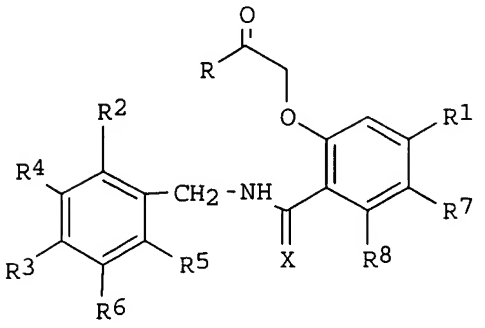
W: AE, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, CA, CH, CN, CR, CU, CZ, DE, DK, DM, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, US, UZ, VN, YU, ZA, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM

RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG

CA 2385798	AA	20010104	CA 2000-2385798	20000623
BR 2000011928	A	20020409	BR 2000-11928	20000623
EP 1198451	A2	20020424	EP 2000-944834	20000623
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, MK, CY, AL				
US 6420426	B1	20020716	US 2000-603817	20000623
JP 2003503381	T2	20030128	JP 2001-506979	20000623
EE 200100708	A	20030217	EE 2001-708	20000623
NZ 516290	A	20040326	NZ 2000-516290	20000623
NO 2001006272	A	20020117	NO 2001-6272	20011220
ZA 2002000300	A	20030613	ZA 2002-300	20020114
BG 106351	A	20020930	BG 2002-106351	20020125
US 2003036558	A1	20030220	US 2002-195964	20020716
PRIORITY APPLN. INFO.:			US 1999-141068P	P 19990625
			US 2000-603817	A3 20000623
			WO 2000-US17377	W 20000623

OTHER SOURCE(S): MARPAT 134:71392

GI



I

AB Disclosed are substituted phenoxyacetic acids [I; R = OH, OCH2CH3, C6H5CH2O, (CH3)2CHCH2CH2O, CH3(CH2)7O, CH3(CH2)3; R1 = H, Cl, F, Br, CH3O, CH3, NO2, CH3S, CF3O, CH3SO2; R2 = F, H; R3 = Br, H, Cl, CH3O, CF3O, CH3; R4 = H, F, CF3, NO2, CH3O, Cl; R5 = H, F; R6 = H, F, CF3, CH3O; R7 = H, CH3, CF3O, NH2, F; R8 = H, F; X = O, S; etc] useful in the treatment of chronic complications arising from diabetes mellitus. Also disclosed are pharmaceutical compns. containing title compds. I, alone or in combination with other therapeutic agents, and methods of treatment employing the compds. and pharmaceutical compns., as well as methods for their synthesis. The pharmaceutical composition contains angiotensin converting enzyme inhibitor(benazepril, captopril, delapril, etc). Thus, title compound I (R1 = Cl; R = OH; X = O; R4 = NO2; R2, R3, R5, R6, R7, R8 each = H) was prepared and tested.

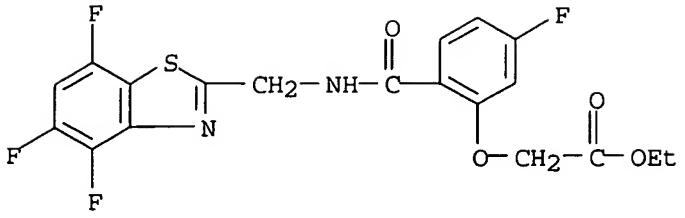
IT **314298-38-9P 314298-39-0P**

RL: RCT (Reactant); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); RACT (Reactant or reagent); USES (Uses)

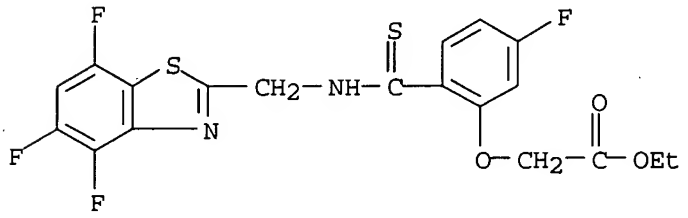
(preparation and effect of substituted phenoxyacetic acids in complications arising from diabetes mellitus)

RN 314298-38-9 CAPLUS

CN Acetic acid, [5-fluoro-2-[[[(4,5,7-trifluoro-2-benzothiazolyl)methyl]amino]carbonyl]phenoxy]-, ethyl ester (9CI) (CA INDEX NAME)

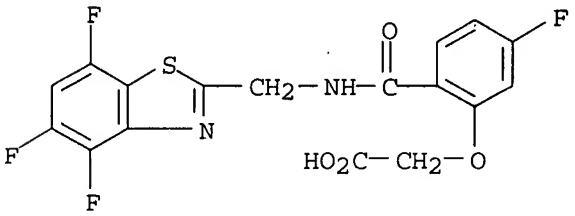


RN 314298-39-0 CAPLUS
CN Acetic acid, [5-fluoro-2-[thioxo[[(4,5,7-trifluoro-2-benzothiazolyl)methyl]amino]methyl]phenoxy]-, ethyl ester (9CI) (CA INDEX NAME)

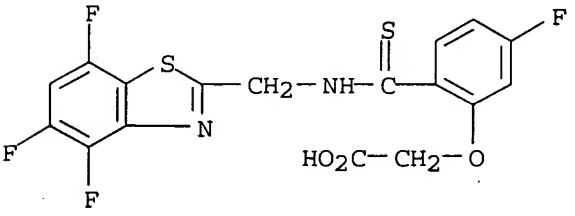


IT **314297-79-5P 314297-80-8P 314297-81-9P 314297-82-0P**
RL: SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(preparation and effect of substituted phenoxyacetic acids in complications arising from diabetes mellitus)

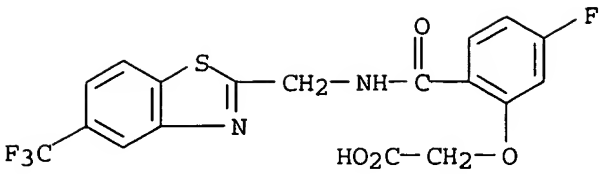
RN 314297-79-5 CAPLUS
CN Acetic acid, [5-fluoro-2-[[[(4,5,7-trifluoro-2-benzothiazolyl)methyl]amino]carbonyl]phenoxy]- (9CI) (CA INDEX NAME)



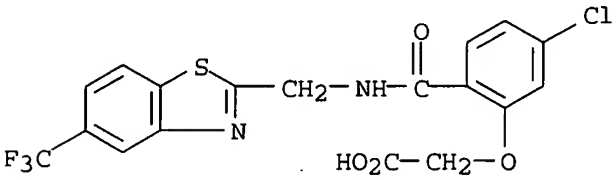
RN 314297-80-8 CAPLUS
CN Acetic acid, [5-fluoro-2-[thioxo[[(4,5,7-trifluoro-2-benzothiazolyl)methyl]amino]methyl]phenoxy]- (9CI) (CA INDEX NAME)



RN 314297-81-9 CAPLUS
CN Acetic acid, [5-fluoro-2-[[[[5-(trifluoromethyl)-2-benzothiazolyl]methyl]amino]carbonyl]phenoxy]- (9CI) (CA INDEX NAME)



RN 314297-82-0 CAPLUS
CN Acetic acid, [5-chloro-2-[[[[5-(trifluoromethyl)-2-benzothiazolyl]methyl]amino]carbonyl]phenoxy]- (9CI) (CA INDEX NAME)



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COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	16.17	346.61

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
CA SUBSCRIBER PRICE	-2.19	-2.19

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